

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
 Data File : BP010022.D
 Acq On : 21 Apr 2022 19:34
 Operator : CG/JU
 Sample : N2373-11
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BFFF0

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
 Title : SVOA CALIBRATION

Signal : TIC: BP010022.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.369	45	48	53	rVB2	11015	13774	0.45%	0.036%
2	3.793	116	120	133	rBV	77135	154541	5.09%	0.399%
3	3.893	133	137	145	rVB2	24848	40542	1.34%	0.105%
4	4.122	173	176	182	rVB5	7150	11369	0.37%	0.029%
5	4.340	210	213	219	rVB8	4595	8780	0.29%	0.023%
6	5.046	328	333	343	rVB4	10898	23104	0.76%	0.060%
7	6.946	652	656	660	rBV6	4760	7359	0.24%	0.019%
8	7.099	675	682	700	rVB	278306	474118	15.62%	1.226%
9	7.263	700	710	722	rBV	192500	323237	10.65%	0.836%
10	7.469	737	745	760	rBV	271694	457615	15.07%	1.183%
11	7.940	818	825	834	rBV	356074	583316	19.22%	1.508%
12	8.187	861	867	871	rBV9	3891	7179	0.24%	0.019%
13	8.646	938	945	956	rBV2	354501	593896	19.56%	1.535%
14	8.763	960	965	971	rVB2	21537	37818	1.25%	0.098%
15	9.098	1014	1022	1033	rBV	261679	452788	14.92%	1.170%
16	9.275	1047	1052	1058	rVB10	3334	6700	0.22%	0.017%
17	9.551	1094	1099	1104	rVB5	12167	22565	0.74%	0.058%
18	9.822	1136	1145	1159	rBV	219055	382506	12.60%	0.989%
19	9.981	1166	1172	1175	rBV8	3458	7101	0.23%	0.018%
20	10.298	1219	1226	1231	rBV	31611	56401	1.86%	0.146%
21	10.363	1231	1237	1251	rVB	398388	673804	22.20%	1.742%
22	10.745	1295	1302	1309	rBV	544060	908503	29.93%	2.349%
23	10.881	1318	1325	1344	rBV	211269	419578	13.82%	1.085%
24	12.181	1539	1546	1549	rBV4	9553	15632	0.51%	0.040%
25	12.269	1554	1561	1568	rVV	80847	140280	4.62%	0.363%
26	12.328	1568	1571	1575	rVV3	8144	14044	0.46%	0.036%
27	12.398	1578	1583	1588	rBV9	4869	9776	0.32%	0.025%
28	13.410	1751	1755	1760	rVV2	13286	19740	0.65%	0.051%
29	13.892	1833	1837	1844	rVB8	5620	10572	0.35%	0.027%
30	13.975	1844	1851	1865	rBV	837522	1213439	39.97%	3.137%
31	14.063	1865	1866	1872	rVB6	4815	7309	0.24%	0.019%
32	14.263	1892	1900	1915	rVV	861334	1318764	43.44%	3.409%
33	14.428	1922	1928	1932	rBV5	7642	12462	0.41%	0.032%
34	14.481	1933	1937	1942	rVB6	5544	8470	0.28%	0.022%
35	14.569	1943	1952	1961	rBV	908775	1375557	45.31%	3.556%
36	14.757	1978	1984	2000	rBV	318881	547289	18.03%	1.415%

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
 Data File : BP010022.D
 Acq On : 21 Apr 2022 19:34
 Operator : CG/JU
 Sample : N2373-11
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BFFF0

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
 Title : SVOA CALIBRATION

37	14.910	2006	2010	2017	rBV7	9019	17896	0.59%	0.046%
38	14.981	2020	2022	2027	rVV6	6876	10813	0.36%	0.028%
39	15.239	2062	2066	2071	rVB8	4948	7247	0.24%	0.019%
40	15.381	2080	2090	2096	rBV10	8346	21900	0.72%	0.057%
41	15.439	2096	2100	2104	rVB5	4991	8642	0.28%	0.022%
42	15.563	2114	2121	2128	rBV	1234455	1777026	58.54%	4.594%
43	15.675	2135	2140	2150	rBV	346511	479770	15.80%	1.240%
44	15.804	2158	2162	2166	rBV3	12980	16646	0.55%	0.043%
45	16.292	2241	2245	2252	rVB10	13125	21268	0.70%	0.055%
46	16.463	2271	2274	2278	rVB5	6703	9340	0.31%	0.024%
47	16.610	2295	2299	2305	rBV5	10422	20162	0.66%	0.052%
48	16.722	2314	2318	2326	rBV9	5734	10494	0.35%	0.027%
49	16.980	2357	2362	2369	rVB7	10104	22764	0.75%	0.059%
50	17.222	2399	2403	2410	rVB4	14927	21870	0.72%	0.057%
51	17.322	2410	2420	2425	rBV	1226857	1761933	58.04%	4.555%
52	17.363	2425	2427	2431	rVV	100734	128780	4.24%	0.333%
53	17.422	2431	2437	2448	rVV	1358332	2031912	66.93%	5.253%
54	17.510	2449	2452	2459	rVB8	13327	23392	0.77%	0.060%
55	17.833	2503	2507	2515	rVV10	10685	25601	0.84%	0.066%
56	17.904	2516	2519	2523	rVB4	10414	12755	0.42%	0.033%
57	17.998	2531	2535	2540	rVB7	14076	19074	0.63%	0.049%
58	18.092	2548	2551	2553	rBV3	7815	11645	0.38%	0.030%
59	18.151	2557	2561	2564	rBV3	41067	63906	2.11%	0.165%
60	18.204	2564	2570	2573	rBV2	130850	183255	6.04%	0.474%
61	18.369	2594	2598	2603	rBV2	47169	82408	2.71%	0.213%
62	18.410	2603	2605	2609	rVB	22885	28456	0.94%	0.074%
63	18.492	2616	2619	2621	rBV4	12457	16652	0.55%	0.043%
64	18.669	2646	2649	2652	rBV2	23224	25503	0.84%	0.066%
65	18.704	2652	2655	2660	rVB6	19870	27494	0.91%	0.071%
66	18.904	2686	2689	2694	rVB6	28400	34221	1.13%	0.088%
67	18.969	2696	2700	2708	rVB6	16206	40613	1.34%	0.105%
68	19.074	2713	2718	2721	rBV3	54639	88380	2.91%	0.228%
69	19.104	2721	2723	2730	rVB4	27422	41861	1.38%	0.108%
70	19.163	2730	2733	2736	rVB2	19740	19088	0.63%	0.049%
71	19.204	2737	2740	2741	rBV2	24430	24954	0.82%	0.065%
72	19.298	2752	2756	2757	rBV	23526	30206	1.00%	0.078%
73	19.327	2757	2761	2767	rVV	279948	366102	12.06%	0.946%
74	19.386	2768	2771	2774	rVB5	14832	17111	0.56%	0.044%
75	19.474	2783	2786	2790	rVB	21235	23612	0.78%	0.061%
76	19.651	2810	2816	2820	rBV	2089145	2850139	93.89%	7.368%

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
 Data File : BP010022.D
 Acq On : 21 Apr 2022 19:34
 Operator : CG/JU
 Sample : N2373-11
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BFFF0

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
 Title : SVOA CALIBRATION

77	19.751	2830	2833	2839	rVB5	28770	46235	1.52%	0.120%
78	19.986	2870	2873	2876	rBV3	30400	49534	1.63%	0.128%
79	20.086	2886	2890	2893	rBV	41035	54156	1.78%	0.140%
80	20.151	2898	2901	2910	rVB7	62653	148203	4.88%	0.383%
81	20.316	2926	2929	2933	rVB2	45858	61287	2.02%	0.158%
82	20.451	2949	2952	2956	rBV2	48239	67576	2.23%	0.175%
83	20.498	2957	2960	2964	rVB2	44242	52773	1.74%	0.136%
84	20.892	3024	3027	3033	rVB	47382	63461	2.09%	0.164%
85	21.110	3059	3064	3072	rBV2	806216	1083675	35.70%	2.801%
86	21.310	3095	3098	3102	rBV	78875	91923	3.03%	0.238%
87	21.404	3107	3114	3118	rVV	1555055	2365097	77.91%	6.114%
88	21.439	3118	3120	3129	rVB	230711	318489	10.49%	0.823%
89	22.045	3217	3223	3232	rVB3	279056	529124	17.43%	1.368%
90	22.533	3302	3306	3321	rVB3	133944	288554	9.51%	0.746%
91	22.792	3345	3350	3362	rVB	563314	867210	28.57%	2.242%
92	23.110	3398	3404	3407	rBV	560778	975640	32.14%	2.522%
93	23.157	3407	3412	3425	rVB	1657315	3035665	100.00%	7.847%
94	23.698	3498	3504	3510	rBV	1086613	2109215	69.48%	5.452%
95	23.857	3524	3531	3539	rBV	881631	1810775	59.65%	4.681%
96	24.115	3571	3575	3584	rVB3	117947	227484	7.49%	0.588%
97	24.509	3635	3642	3649	rVB	556068	1202305	39.61%	3.108%
98	24.598	3651	3657	3669	rVV	346429	784525	25.84%	2.028%
99	26.409	3958	3965	3981	rVB2	336368	1054846	34.75%	2.727%
100	27.386	4126	4131	4142	rVB6	195378	611667	20.15%	1.581%

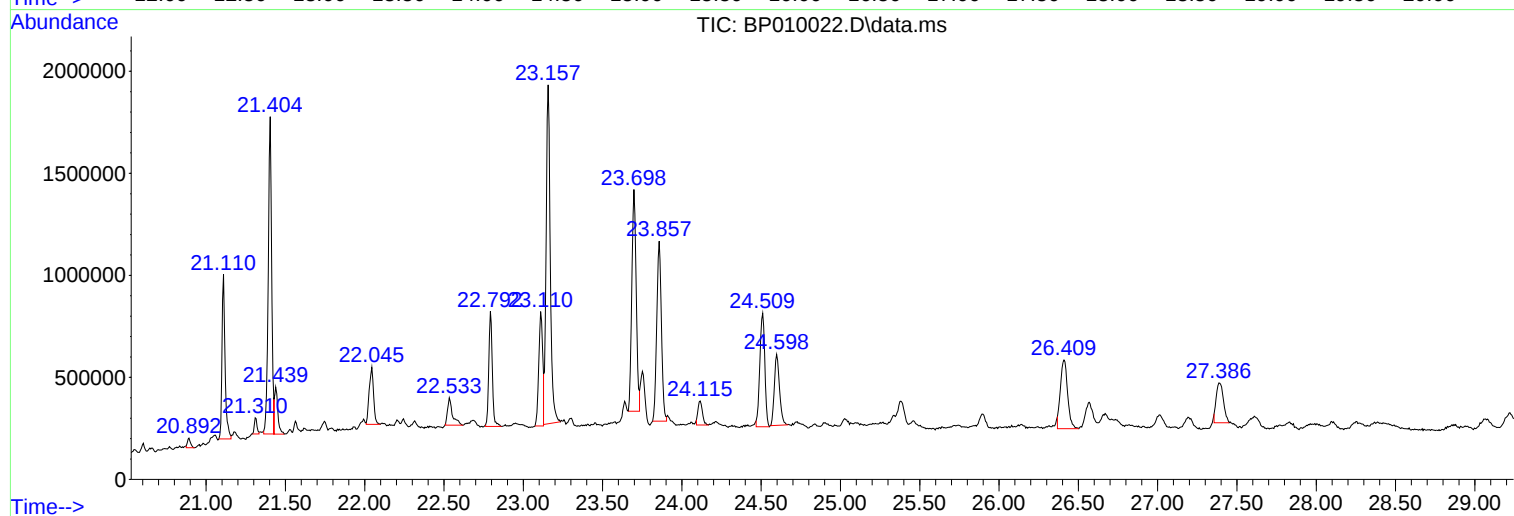
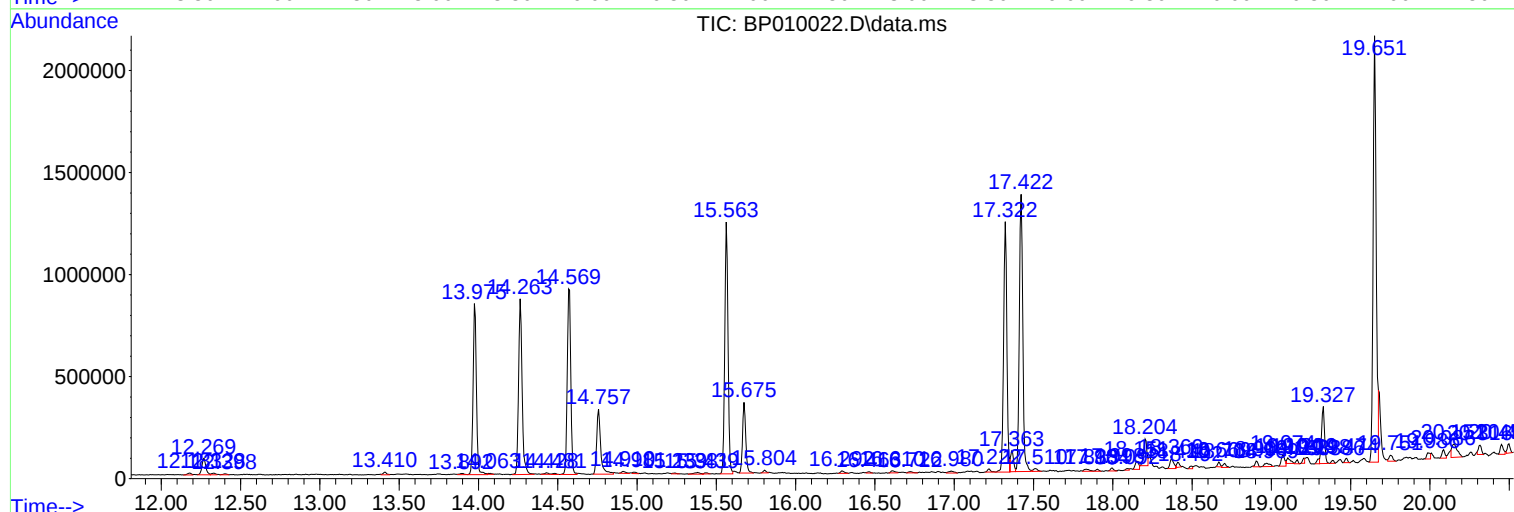
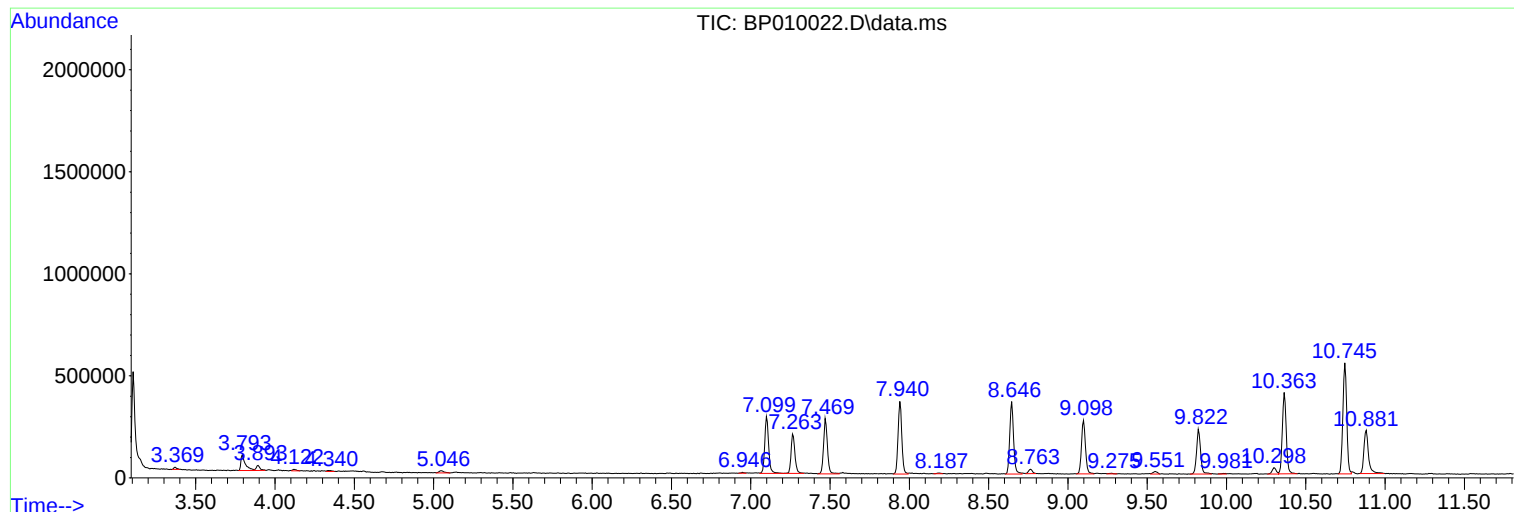
Sum of corrected areas: 38684238

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010022.D
Acq On : 21 Apr 2022 19:34
Operator : CG/JU
Sample : N2373-11
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BFFF0

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010022.D
Acq On : 21 Apr 2022 19:34
Operator : CG/JU
Sample : N2373-11
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BFFF0

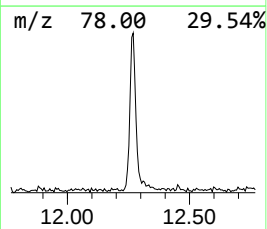
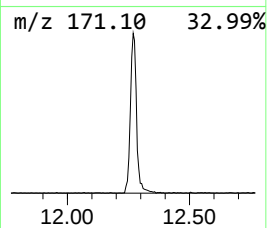
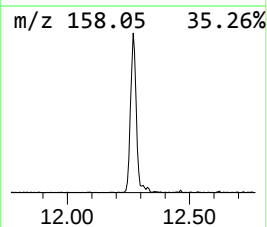
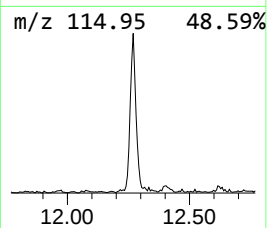
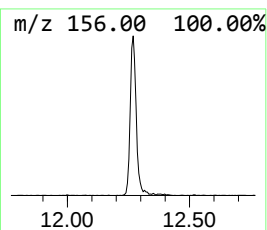
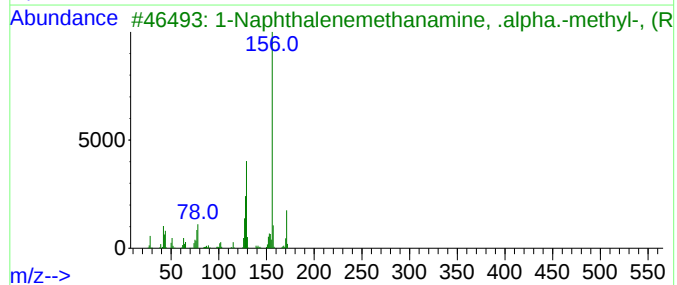
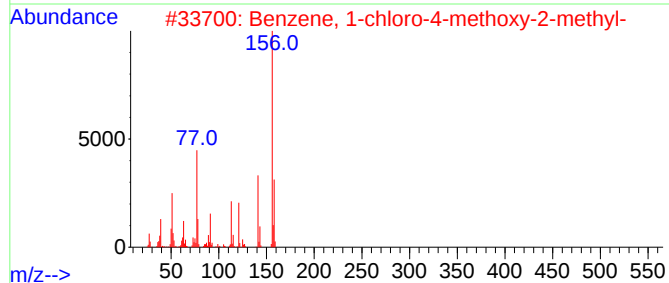
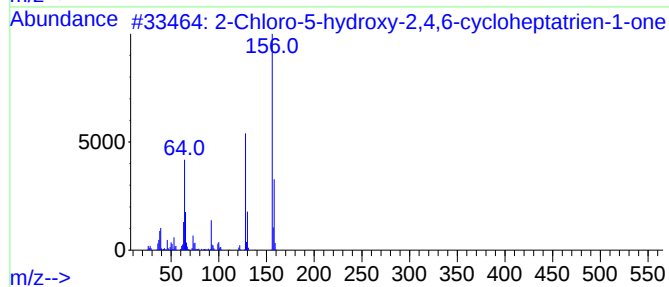
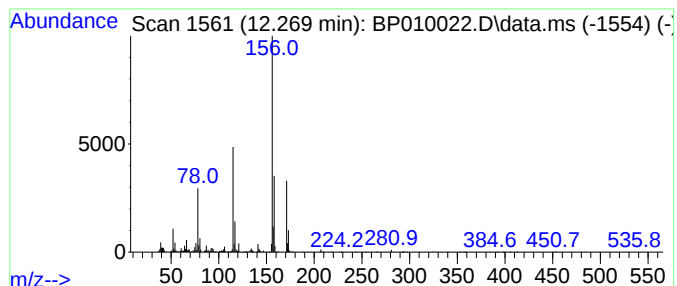
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 unknown-01 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.269	3.09 ng/ul	140280	Naphthalene-d8	10.746

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Chloro-5-hydroxy-2,4,6-cyclohe...	156	C7H5ClO2	007009-16-7	43
2			Benzene, 1-chloro-4-methoxy-2-me...	156	C8H9ClO	013334-71-9	43
3			1-Naphthalenemethanamine, .alpha...	171	C12H13N	003886-70-2	43
4			6-Isopropylquinoline	171	C12H13N	000135-79-5	43
5			1,2,3,3a-Tetrahydropentalene, 1,...	171	C12H13N	1000217-17-0	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010022.D
Acq On : 21 Apr 2022 19:34
Operator : CG/JU
Sample : N2373-11
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BFFF0

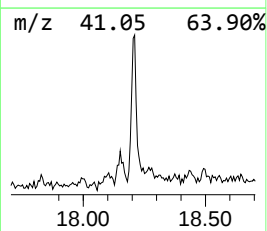
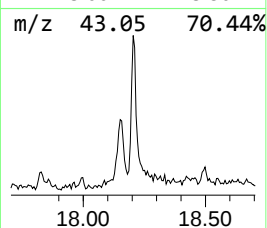
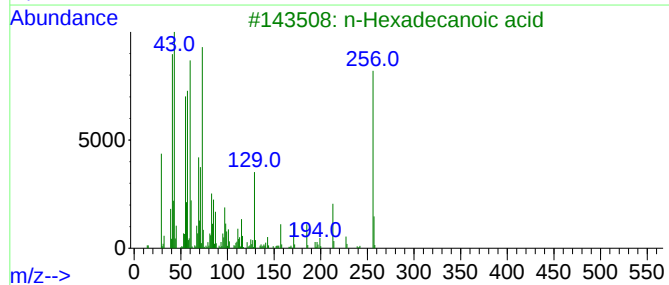
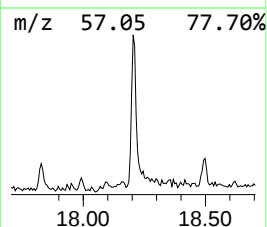
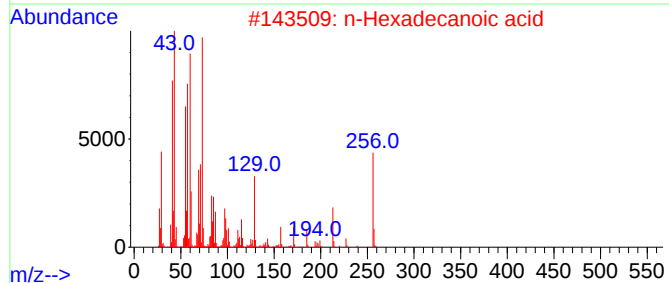
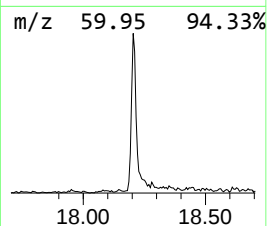
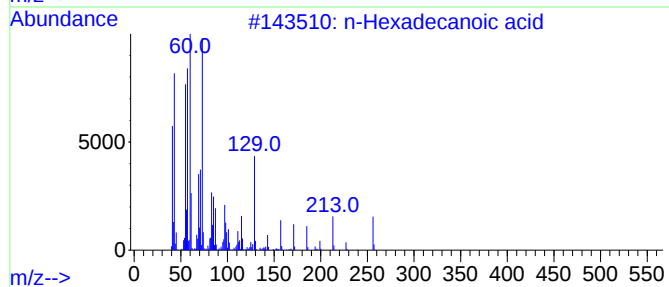
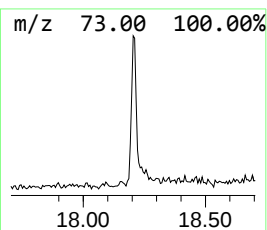
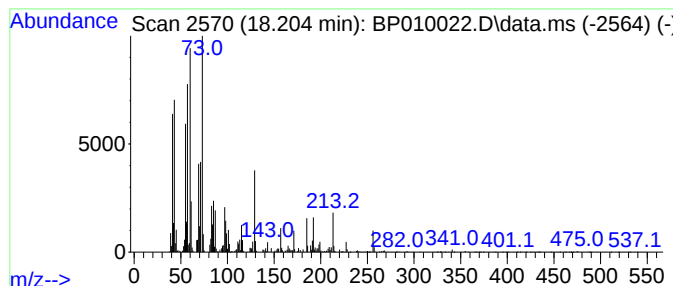
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 n-Hexadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.204	2.08 ng/ul	183255	Phenanthrene-d10	17.322

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	93
5			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010022.D
Acq On : 21 Apr 2022 19:34
Operator : CG/JU
Sample : N2373-11
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BFFF0

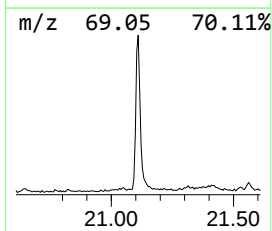
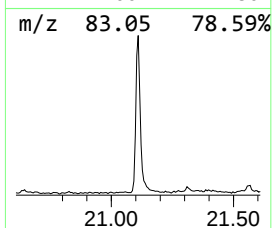
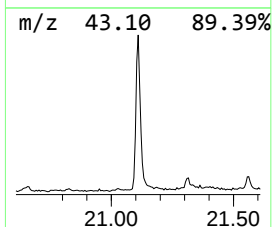
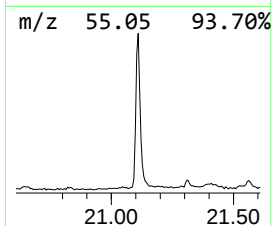
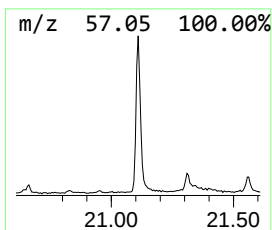
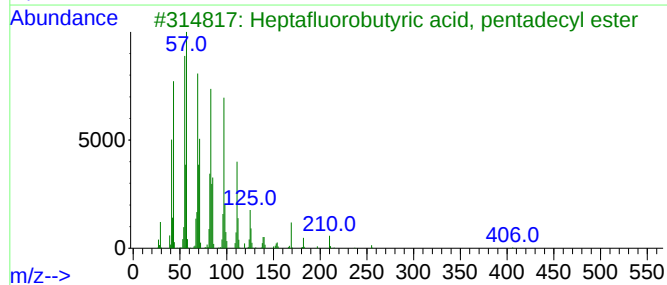
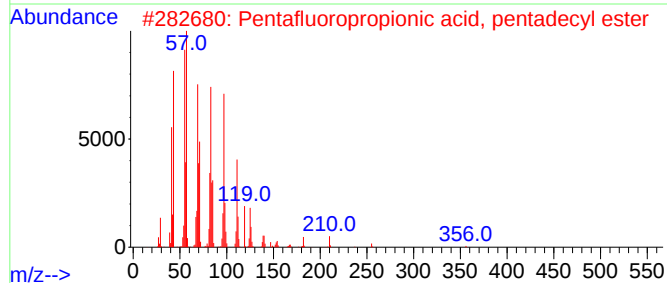
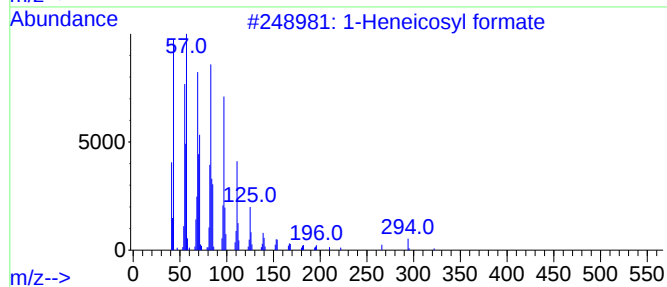
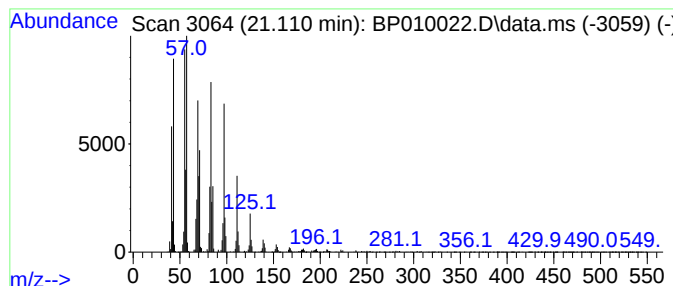
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 1-Heneicosyl formate Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.110	9.16 ng/ul	1083680	Chrysene-d12	21.404

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Heneicosyl formate	340	C22H44O2	077899-03-7	96
2			Pentafluoropropionic acid, penta...	374	C18H31F5O2	959092-08-1	94
3			Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	94
4			Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	94
5			1-Hexacosene	364	C26H52	018835-33-1	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010022.D
Acq On : 21 Apr 2022 19:34
Operator : CG/JU
Sample : N2373-11
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BFFF0

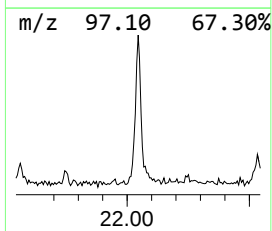
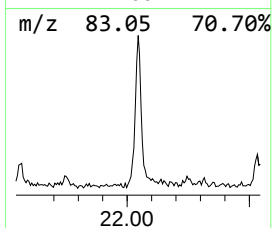
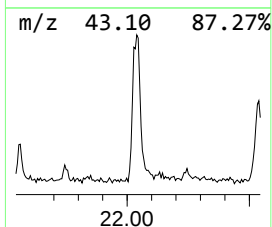
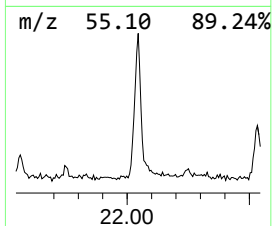
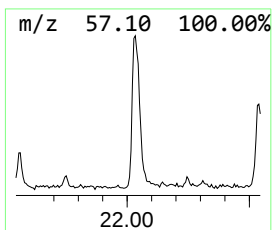
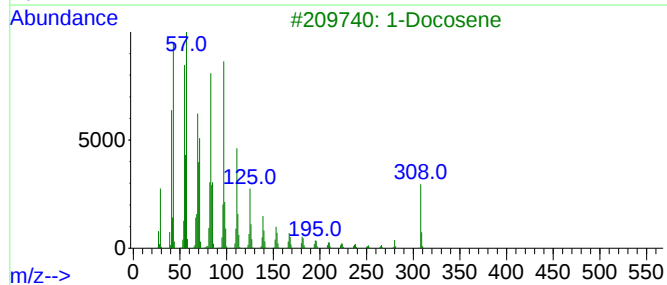
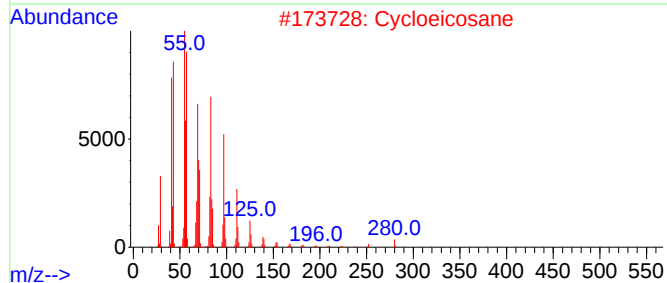
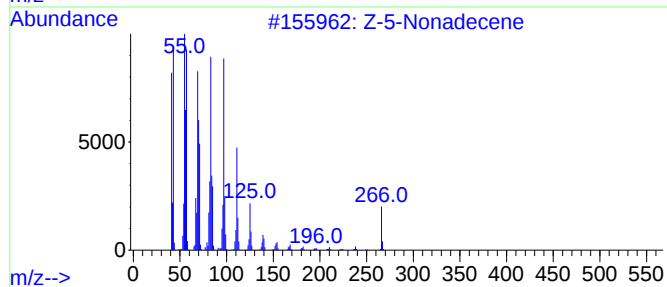
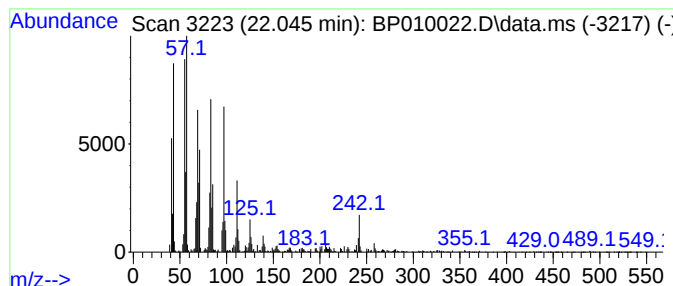
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.045	4.47 ng/ul	529124	Chrysene-d12	21.404

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Z-5-Nonadecene	266	C19H38	1000131-11-8	97
2		Cycloeicosane	280	C20H40	000296-56-0	95
3		1-Docosene	308	C22H44	001599-67-3	95
4		1-Docosene	308	C22H44	001599-67-3	93
5		9-Tricosene, (Z)-	322	C23H46	027519-02-4	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010022.D
Acq On : 21 Apr 2022 19:34
Operator : CG/JU
Sample : N2373-11
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BFFF0

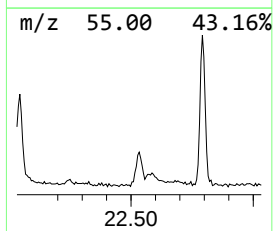
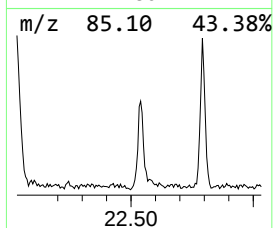
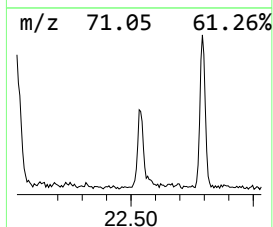
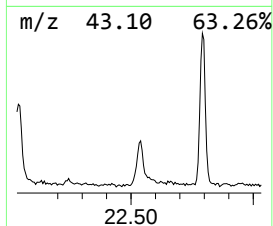
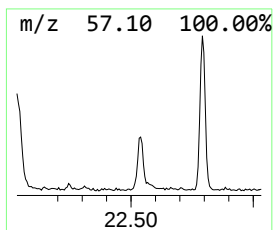
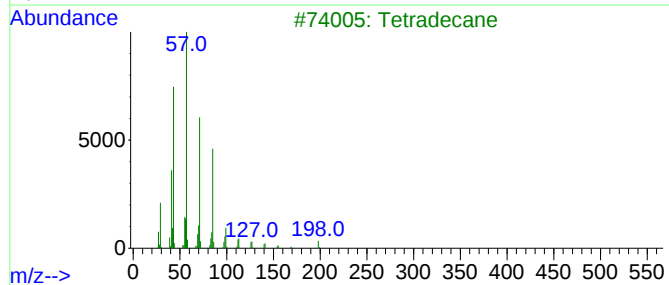
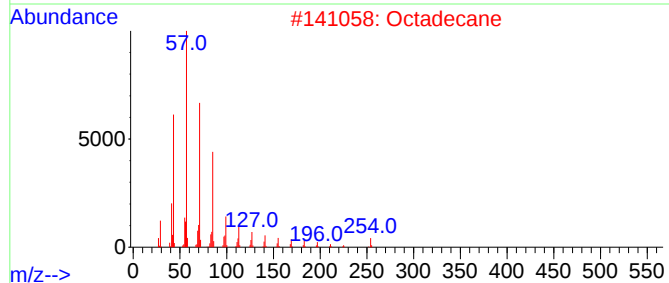
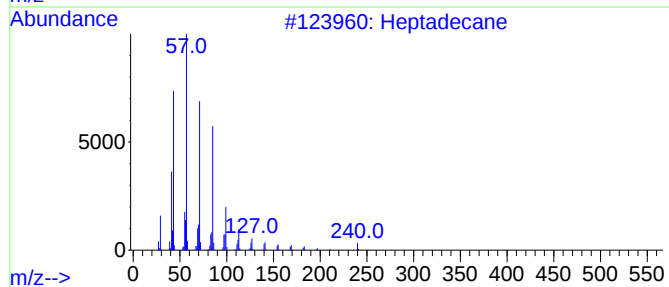
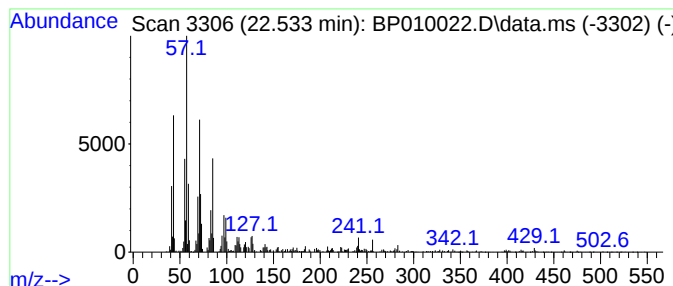
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.533	2.44 ng/ul	288554	Chrysene-d12	21.404

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecane	240	C17H36	000629-78-7	92
2		Octadecane	254	C18H38	000593-45-3	91
3		Tetradecane	198	C14H30	000629-59-4	90
4		Heptadecane	240	C17H36	000629-78-7	90
5		Eicosane	282	C20H42	000112-95-8	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010022.D
Acq On : 21 Apr 2022 19:34
Operator : CG/JU
Sample : N2373-11
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BFFF0

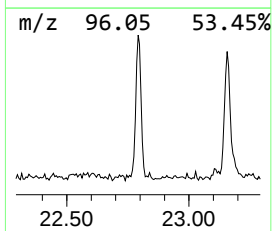
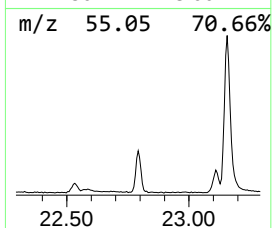
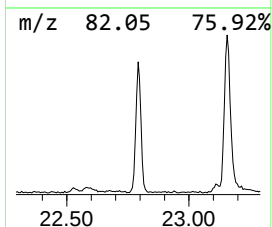
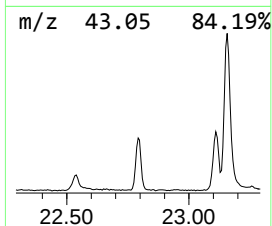
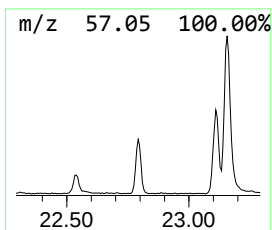
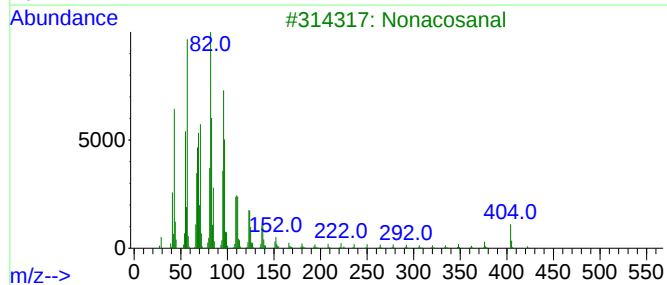
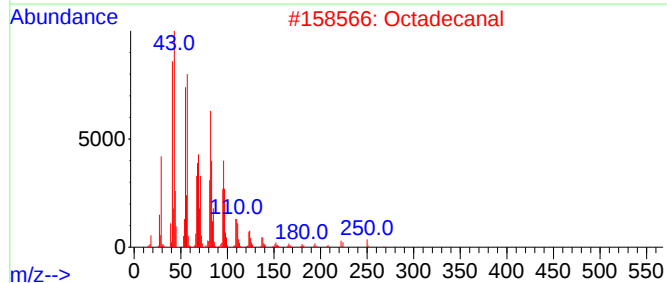
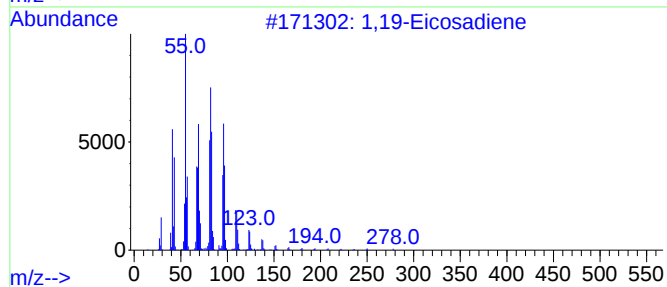
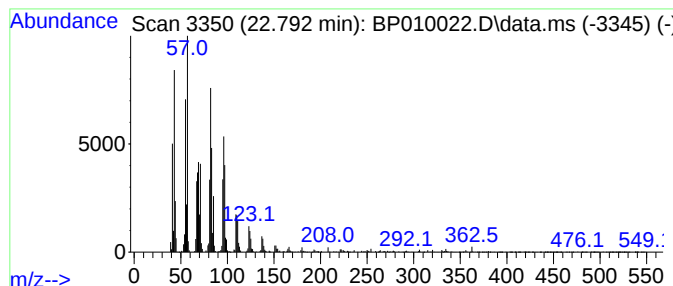
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 1,19-Eicosadiene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.792	9.58 ng/ul	867210	Perylene-d12	23.857

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,19-Eicosadiene			278	C20H38	014811-95-1	95
2	Octadecanal			268	C18H36O	000638-66-4	94
3	Nonacosanal			422	C29H58O	072934-04-4	94
4	Docosanal			324	C22H44O	057402-36-5	91
5	Oxirane, hexadecyl-			268	C18H36O	007390-81-0	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010022.D
Acq On : 21 Apr 2022 19:34
Operator : CG/JU
Sample : N2373-11
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BFFF0

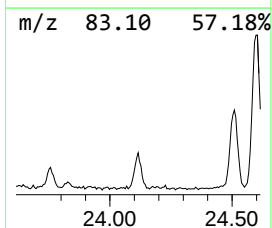
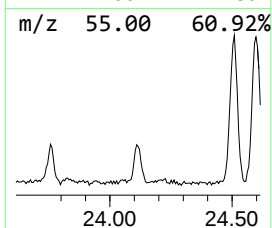
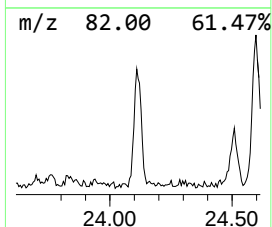
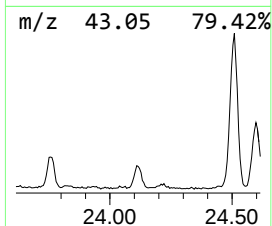
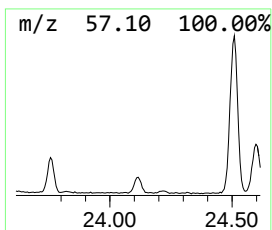
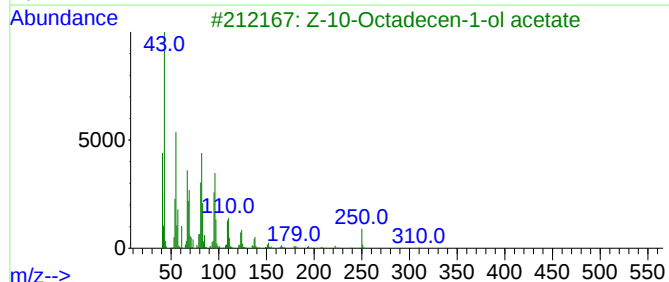
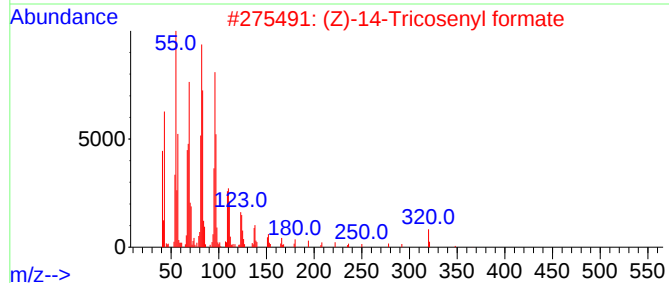
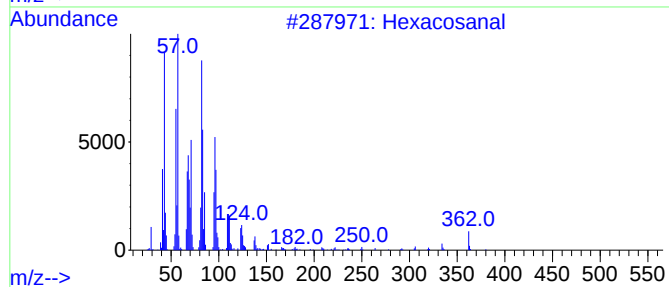
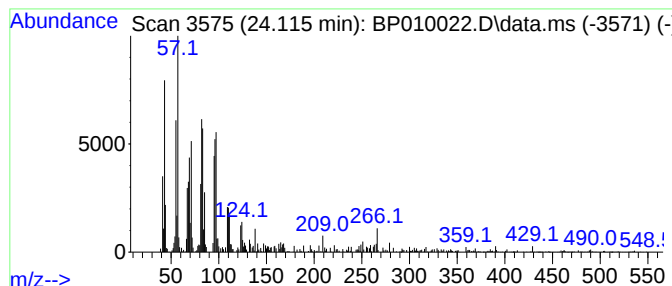
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Hexacosanal Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.115	2.51 ng/ul	227484	Perylene-d12	23.857

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexacosanal	380	C26H52O	026627-85-0	83
2			(Z)-14-Tricosenyl formate	366	C24H46O2	077899-10-6	83
3			Z-10-Octadecen-1-ol acetate	310	C20H38O2	1000131-07-2	78
4			Dotriacontanal	464	C32H64O	057878-00-9	76
5			Tridecanal	198	C13H26O	010486-19-8	76



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010022.D
Acq On : 21 Apr 2022 19:34
Operator : CG/JU
Sample : N2373-11
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BFFF0

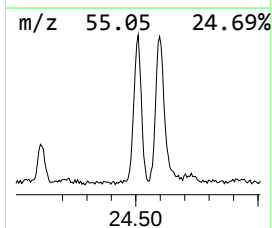
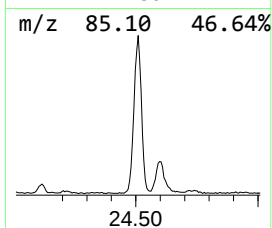
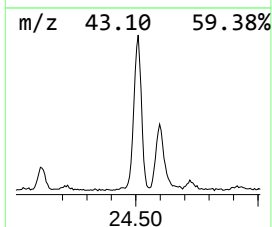
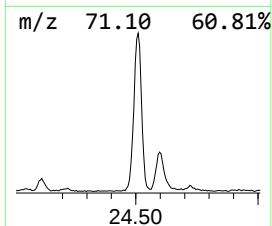
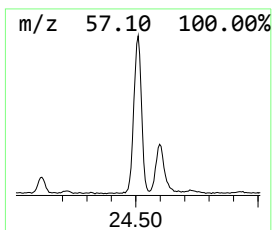
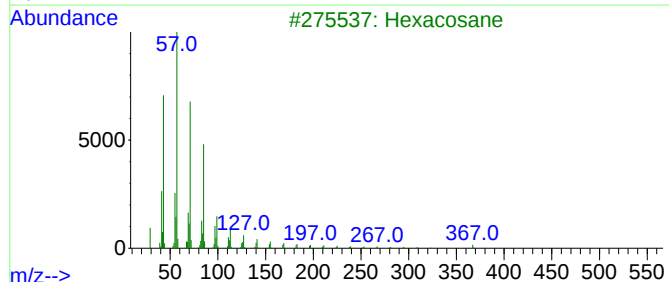
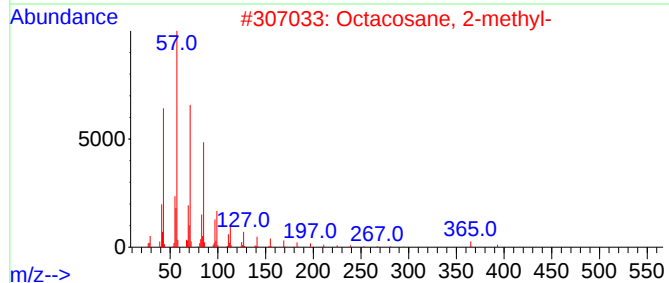
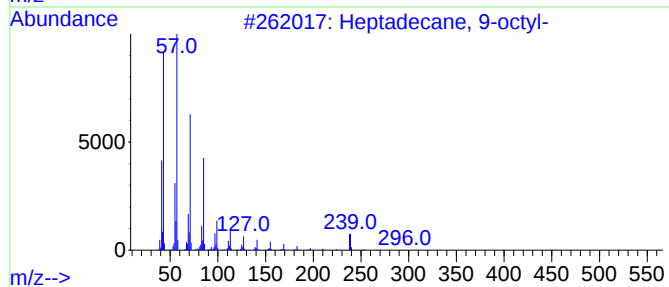
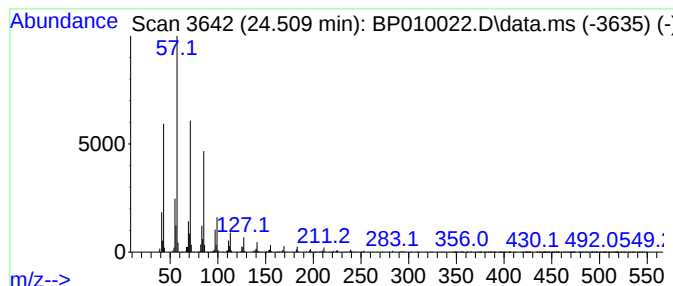
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.509	13.28 ng/ul	1202310	Perylene-d12	23.857

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	93
2		Octacosane, 2-methyl-	408	C29H60	001560-98-1	91
3		Hexacosane	366	C26H54	000630-01-3	91
4		Tetratetracontane	619	C44H90	007098-22-8	91
5		Octacosane	394	C28H58	000630-02-4	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010022.D
Acq On : 21 Apr 2022 19:34
Operator : CG/JU
Sample : N2373-11
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BFFF0

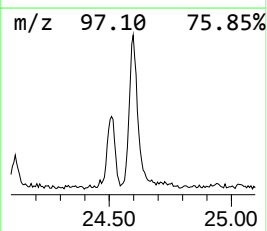
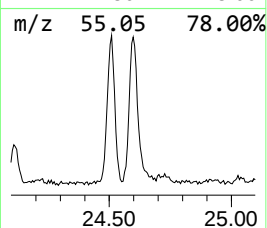
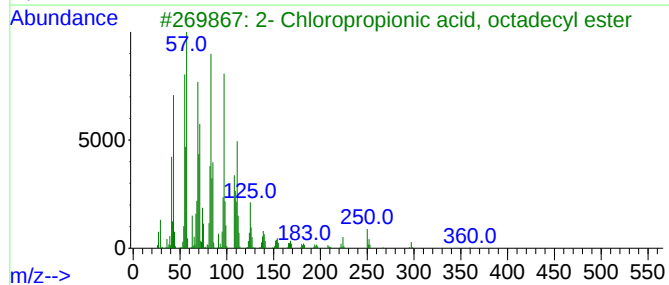
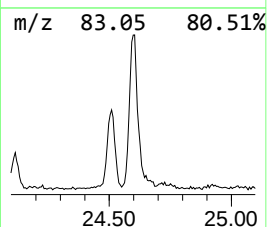
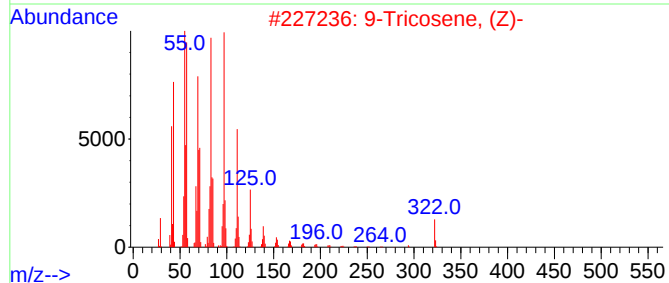
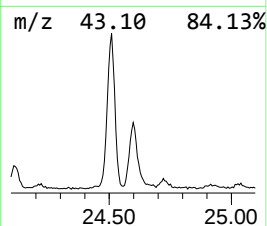
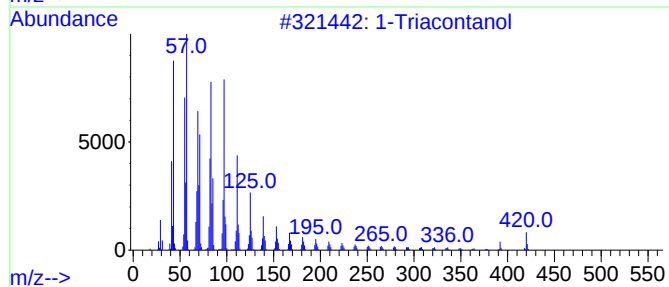
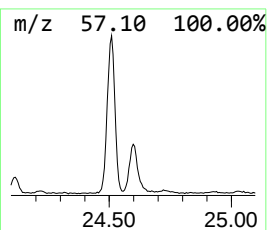
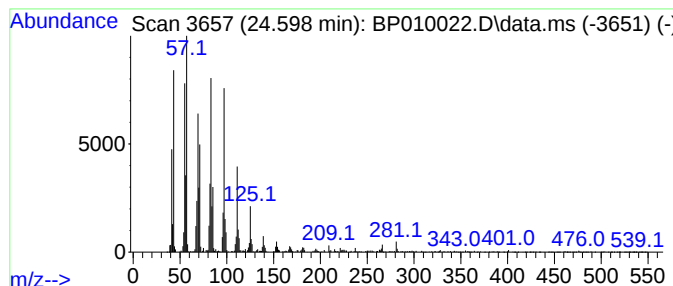
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 1-Triacontanol Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.598	8.67 ng/ul	784525	Perylene-d12	23.857

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Triacontanol			438	C30H62O	000593-50-0	96
2	9-Tricosene, (Z)-			322	C23H46	027519-02-4	95
3	2- Chloropropionic acid, octadec...			360	C21H41ClO2	088104-31-8	94
4	Nonacos-1-ene			406	C29H58	018835-35-3	94
5	1-Heptacosanol			396	C27H56O	002004-39-9	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010022.D
Acq On : 21 Apr 2022 19:34
Operator : CG/JU
Sample : N2373-11
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BFFF0

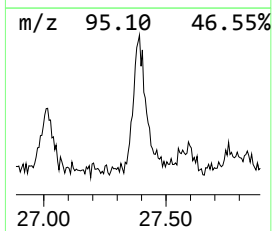
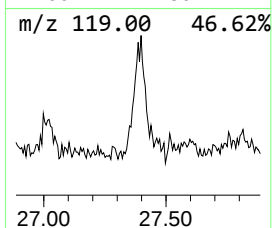
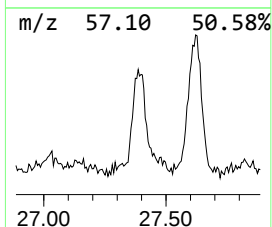
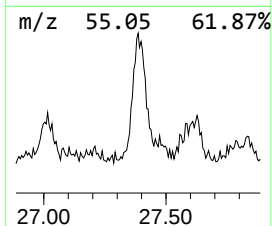
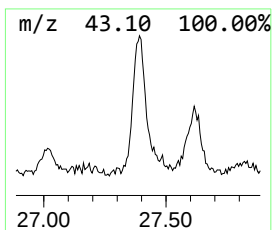
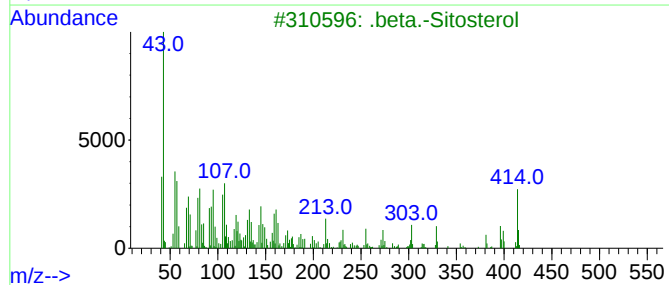
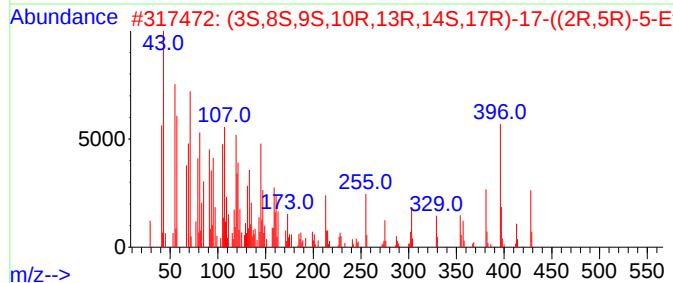
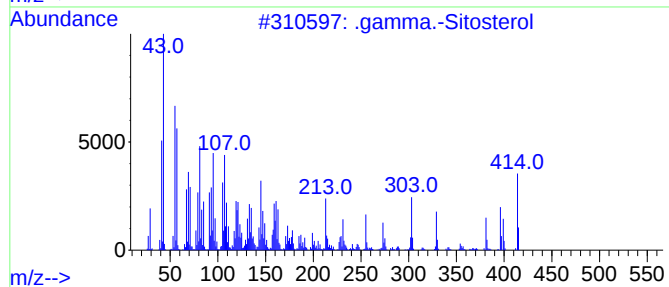
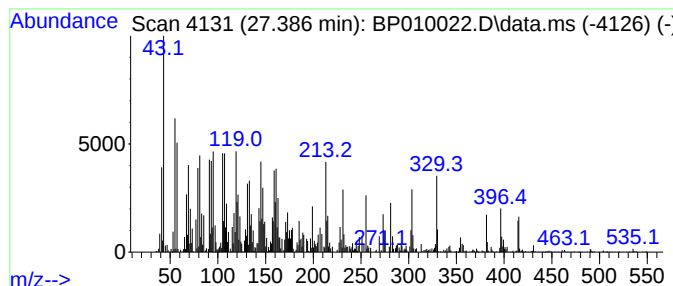
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 unknown-02 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
27.386	6.76 ng/ul	611667	Perylene-d12	23.857

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	.	gamma.-	Sitosterol	414	C29H50O	000083-47-6	47
2	(3S,8S,9S,10R,13R,14S,17R)-17-((...			428	C30H52O	020194-50-7	45
3	.	beta.-	Sitosterol	414	C29H50O	000083-46-5	30
4	.	beta.-	Sitosterol	414	C29H50O	000083-46-5	17
5	(3S,8S,9S,10R,13R,14S,17R)-17-((...			414	C29H50O	029365-29-5	16



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010022.D
Acq On : 21 Apr 2022 19:34
Operator : CG/JU
Sample : N2373-11
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BFFF0

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown-01	12.269	3.1	ng/ul	140280	2	10.746	908503	20.0
n-Hexadecanoic ...	18.204	2.1	ng/ul	183255	4	17.322	1761930	20.0
1-Heneicosyl fo...	21.110	9.2	ng/ul	1083680	5	21.404	2365100	20.0
(DEL) Alkane: S...	22.045	4.5	ng/ul	529124	5	21.404	2365100	20.0
(DEL) Alkane: S...	22.533	2.4	ng/ul	288554	5	21.404	2365100	20.0
1,19-Eicosadiene	22.792	9.6	ng/ul	867210	6	23.857	1810780	20.0
Hexacosanal	24.115	2.5	ng/ul	227484	6	23.857	1810780	20.0
(DEL) Alkane: S...	24.509	13.3	ng/ul	1202310	6	23.857	1810780	20.0
1-Triacontanol	24.598	8.7	ng/ul	784525	6	23.857	1810780	20.0
unknown-02	27.386	6.8	ng/ul	611667	6	23.857	1810780	20.0